

09/438,365

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1204BXD

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * Welcome to STN International * * * * *

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
NEWS 2 "Ask CAS" for self-help around the clock
NEWS 3 FEB 25 CA/CAPLUS - Russian Agency for Patents and Trademarks
(ROSPATENT) added to list of core patent offices covered
NEWS 4 FEB 28 PATDPAFULL - New display fields provide for legal status
data from INPADOC
NEWS 5 FEB 28 BABS - Current-awareness alerts (SDIs) available
NEWS 6 FEB 28 MEDLINE/LMEDLINE reloaded
NEWS 7 MAR 02 GBFULL: New full-text patent database on STN
NEWS 8 MAR 03 REGISTRY/ZREGISTRY - Sequence annotations enhanced
NEWS 9 MAR 03 MEDLINE file segment of TOXCENTER reloaded
NEWS 10 MAR 22 KOREAPAT now updated monthly; patent information enhanced
NEWS 11 MAR 22 Original IDE display format returns to REGISTRY/ZREGISTRY
NEWS 12 MAR 22 PATDPASPC - New patent database available
NEWS 13 MAR 22 REGISTRY/ZREGISTRY enhanced with experimental property tags
NEWS 14 APR 04 EPFULL enhanced with additional patent information and new
fields
NEWS 15 APR 04 EMBASE - Database reloaded and enhanced
NEWS 16 APR 18 New CAS Information Use Policies available online

NEWS EXPRESS JANUARY 10 CURRENT WINDOWS VERSION IS V7.01a, CURRENT
MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
AND CURRENT DISCOVER FILE IS DATED 10 JANUARY 2005

NEWS HOURS STN Operating Hours Plus Help Desk Availability
NEWS INTER General Internet Information
NEWS LOGIN Welcome Banner and News Items
NEWS PHONE Direct Dial and Telecommunication Network Access to STN
NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that
specific topic.

All use of STN is subject to the provisions of the STN Customer
agreement. Please note that this agreement limits use to scientific
research. Use for software development or design or implementation
of commercial gateways or other similar uses is prohibited and may
result in loss of user privileges and other penalties.

* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 14:48:27 ON 25 APR 2005

=> fil reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

FULL ESTIMATED COST

ENTRY	SESSION
0.21	0.21

FILE 'REGISTRY' ENTERED AT 14:48:38 ON 25 APR 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2005 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.

STRUCTURE FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9
DICTIONARY FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

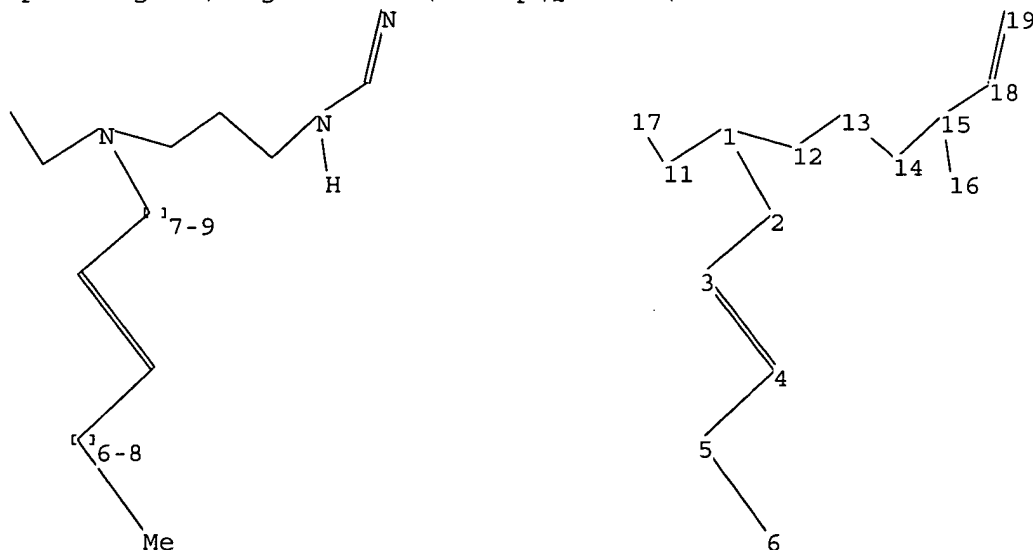
*
* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added, *
* effective March 20, 2005. A new display format, IDERL, is now *
* available and contains the CA role and document type information. *
*

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
information enter HELP PROP at an arrow prompt in the file or refer
to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>

Uploading C:\Program Files\Stnexp\Queries\09438365.str



chain nodes :

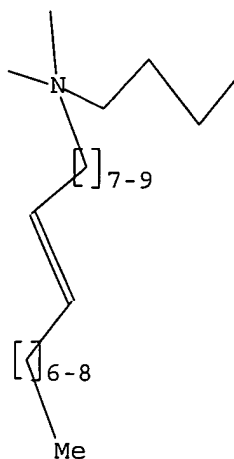
1 2 3 4 5 6 11 12 13 14 15 16 17 18 19
chain bonds :
1-2 1-11 1-12 2-3 3-4 4-5 5-6 11-17 12-13 13-14 14-15 15-16 15-18
18-19
exact/norm bonds :
1-2 1-11 1-12 14-15 15-18 18-19
exact bonds :
2-3 3-4 4-5 5-6 11-17 12-13 13-14 15-16

Match level :
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 11:CLASS 12:CLASS 13:CLASS
14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS

L1 STRUCTURE UPLOADED

=> d query

L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> s l1

SAMPLE SEARCH INITIATED 14:48:57 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 292 TO ITERATE

100.0% PROCESSED 292 ITERATIONS
SEARCH TIME: 00.00.01

10 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 4815 TO 6865
PROJECTED ANSWERS: 11 TO 389

L2 10 SEA SSS SAM L1

=> s l1 full

FULL SEARCH INITIATED 14:49:03 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 5965 TO ITERATE

100.0% PROCESSED 5965 ITERATIONS
SEARCH TIME: 00.00.01

158 ANSWERS

L3 158 SEA SSS FUL L1

=> fil caplus

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

161.33

161.54

FILE 'CAPLUS' ENTERED AT 14:49:11 ON 25 APR 2005

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 25 Apr 2005 VOL 142 ISS 18

FILE LAST UPDATED: 24 Apr 2005 (20050424/ED)

New CAS Information Use Policies, enter HELP USAGETERMS for details.

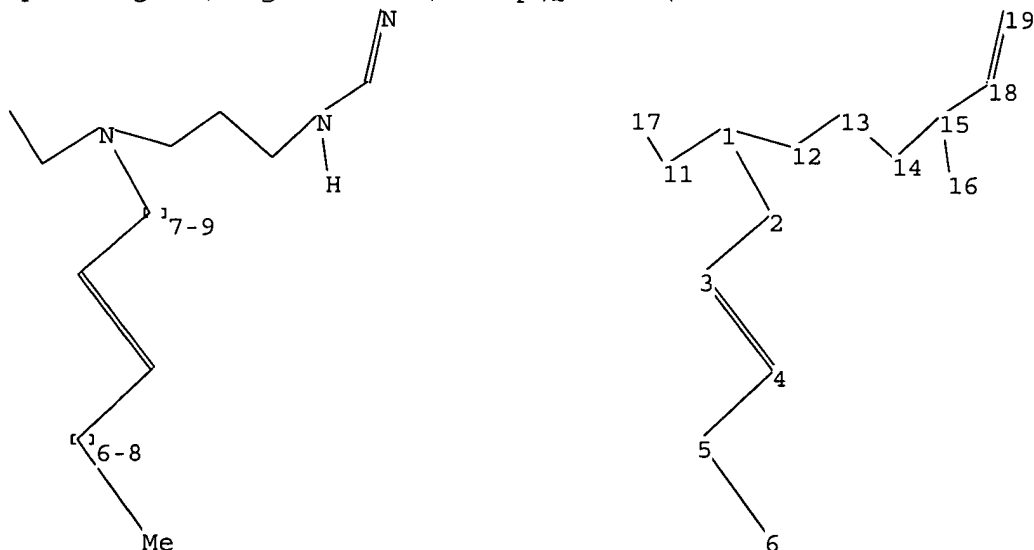
This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l3

L4 281 L3

=>

Uploading C:\Program Files\Stnexp\Queries\09438365.str



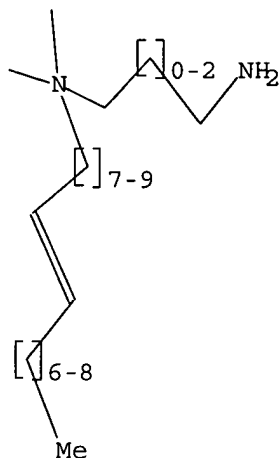
chain nodes :

1 2 3 4 5 6 11 12 13 14 15 16 17 18 19
chain bonds :
1-2 1-11 1-12 2-3 3-4 4-5 5-6 11-17 12-13 13-14 14-15 15-16 15-18
18-19
exact/norm bonds :
1-2 1-11 1-12 14-15 15-18 18-19
exact bonds :
2-3 3-4 4-5 5-6 11-17 12-13 13-14 15-16

Match level :
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 11:CLASS 12:CLASS 13:CLASS
14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS

L5 STRUCTURE UPLOADED

=> d query
L5 STR



Structure attributes must be viewed using STN Express query preparation.

=> s 15

REGISTRY INITIATED

Substance data SEARCH and crossover from CAS REGISTRY in progress...
Use DISPLAY HITSTR (or FHITSTR) to directly view retrieved structures.

SAMPLE SEARCH INITIATED 14:50:20 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 99 TO ITERATE

100.0% PROCESSED 99 ITERATIONS 1 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 1384 TO 2576
PROJECTED ANSWERS: 1 TO 80

L6 1 SEA SSS SAM L5

L7 0 L6

=> fil reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.45	163.32

FILE 'REGISTRY' ENTERED AT 14:50:32 ON 25 APR 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2005 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.

STRUCTURE FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9
DICTIONARY FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

*
* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added, *
* effective March 20, 2005. A new display format, IDERL, is now *
* available and contains the CA role and document type information. *
*

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
information enter HELP PROP at an arrow prompt in the file or refer
to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>

=> s l5 full

FULL SEARCH INITIATED 14:50:36 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 1967 TO ITERATE

100.0% PROCESSED 1967 ITERATIONS 6 ANSWERS
SEARCH TIME: 00.00.01

L8 6 SEA SSS FUL L5

=> fil caplus

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
----------------------	------------	-------

	ENTRY	SESSION
FULL ESTIMATED COST	161.33	324.65

FILE 'CAPLUS' ENTERED AT 14:50:41 ON 25 APR 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 25 Apr 2005 VOL 142 ISS 18
FILE LAST UPDATED: 24 Apr 2005 (20050424/ED)

New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 18

L9 2 L8

=> d 19 1-2 abs ibib hitstr

L9 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2005 ACS on STN

AB The title compds. are useful as emulsifying agents and as fat-liquoring or softening agents for leather, paper, and textiles. Inter-mediate RNHCH₂CH(OH)CONH₂ (I) and RN(CH₂CH(OH)-CONH₂)₂ (II) were prepared by reacting 1 or 2 moles glycidamide with 1 mole of amine RNH₂ in 25 aqueous,

aqueous methanolic, or methanolic solution at 50-60° during 1-2 hr. Quaternization of I and II was carried out by treating with an appropriate halide in H₂O, aqueous methanolic or methanolic solution at 30-50° for 2-3 hr then at 50° for 2-3 hr. Aqueous solns. containing 0.3 g/l. [C₁₈H₃₉-NMe(CH₂CH(OH)CONH₂)₂]+MeSO₄- were used to treat cotton and polycaprolactam fabrics giving material with a soft pleasant hand and hydrophilic in character.

ACCESSION NUMBER: 1971:87345 CAPLUS
DOCUMENT NUMBER: 74:87345
TITLE: (β-Carbamyl-β-hydroxyethyl)alkylammonium salts
PATENT ASSIGNEE(S): Badische Anilin- & Soda-Fabrik AG
SOURCE: Brit., 12 pp.
CODEN: BRXOAA
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
GB 1211040		19701104	GB	
FR 1592740			FR	
US 3632623		19720000	US	
			DE	19671121

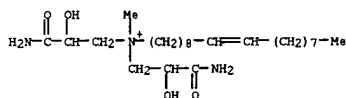
PRIORITY APPLN. INFO.:

IT 32671-74-2P
RI: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)

RN 32671-74-2 CAPLUS
CN Ammonium, bis(2-carbamyl-2-hydroxyethyl)methyl-9-octadecenyl-, methyl sulfate, (2)- (8CI) (CA INDEX NAME)

CM 1

CRN 48077-37-8
CMF C25 H50 N3 O4



CM 2

CRN 21228-90-0
CMF C H3 O4 S

L9 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2005 ACS on STN

AB The title compds. [RR₁R₂NCH₂CONH₂]+X- (I), useful as antistatic agents for synthetic fibers, were prepared by reaction of a tertiary amine, RR₁R₂N (II) with XCH₂CONH₂ (III, X = halogen). Thus, a solution of 297 II (R = n-C₁₈H₃₇,

R₁ = R₂ = Me) and 93.5 III (X = Cl) in 390 (weight parts) MeOH was refluxed

4 hrs., evaporated in vacuo at 50-60°, cooled, and filtered to give I (R = n-C₁₈H₃₇, R₁ = R₂ = Me, X = Cl), straw-colored solids. Other I prepared were (R, R₁, R₂, X): stearamidopropyl, CH₂CH₂OH, CH₂CH₂OH, Cl; n-C₁₁H₂₃CO₂CH₂CH₂, CH₂CH₂OH, CH₂CH₂OH, Cl; oleyl, CH₂CH₂OH, CH₂CH₂OH, Cl; Me n-C₁₁H₂₃CONH-CH₂CH₂, C₁₁H₂₃CONHCH₂CH₂, Br; and Me, oleyl, oleyl, Br; lauryl, (NR₁R₂) morpholino, Cl; oleyl, (CH₂CH₂)₈Al, (CH₂CH₂)₈Al, Br.

ACCESSION NUMBER: 1969:114596 CAPLUS
DOCUMENT NUMBER: 70:114596
TITLE: Quaternary ammonium salts
INVENTOR(S): Fujimoto, Takehiko; Saito, Toshio; Suwada, Ataru; Ohno, Satoyoshi
PATENT ASSIGNEE(S): Sanyo Chemical Industry Co., Ltd.
SOURCE: Jpn. Tokkyo Koho, 3 pp.
CODEN: JAXXAD
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

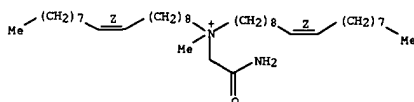
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 43013966	B4	19680613	JP	19640616

IT 23208-83-5P

RI: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)

RN 23208-83-5 CAPLUS
CN Ammonium, (carbamylmethyl)methyl-9-octadecenyl-, bromide, (Z,Z)- (8CI)
(CA INDEX NAME)

Double bond geometry as shown.



• Br⁻

L9 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2005 ACS on STN

(Continued)

Me-O-SO₃⁻


```
=> fil reg
COST IN U.S. DOLLARS          SINCE FILE      TOTAL
                                ENTRY      SESSION
FULL ESTIMATED COST          11.23      335.88

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)  SINCE FILE      TOTAL
                                                ENTRY      SESSION
CA SUBSCRIBER PRICE          -1.46      -1.46
```

FILE 'REGISTRY' ENTERED AT 14:52:11 ON 25 APR 2005
 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
 PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
 COPYRIGHT (C) 2005 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
 provided by InfoChem.

STRUCTURE FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9
 DICTIONARY FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

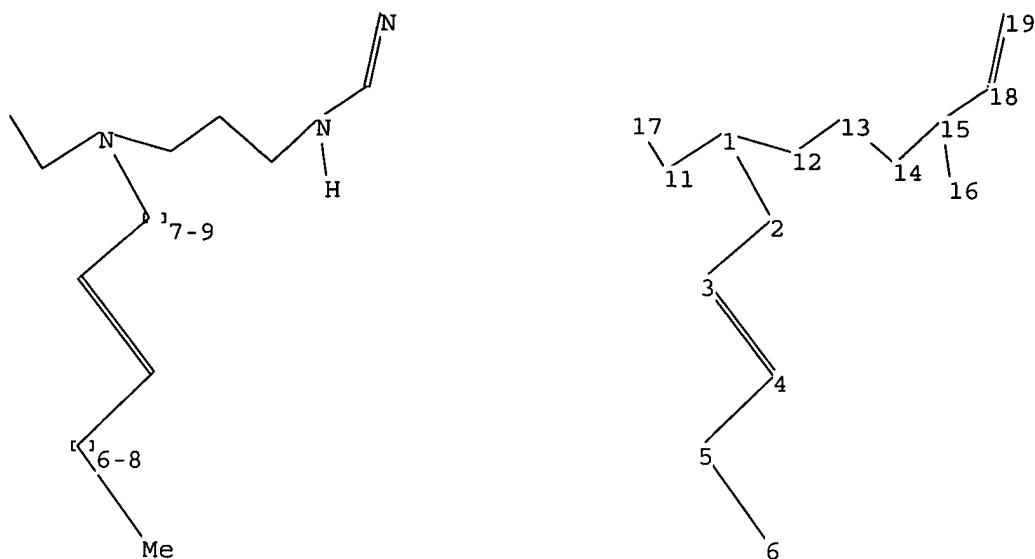
Please note that search-term pricing does apply when
 conducting SmartSELECT searches.

```
*****
*
* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added,  *
* effective March 20, 2005. A new display format, IDERL, is now    *
* available and contains the CA role and document type information. *
*
*****
```

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
 information enter HELP PROP at an arrow prompt in the file or refer
 to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

```
=>
Uploading C:\Program Files\Stnexp\Queries\09438365.str
```



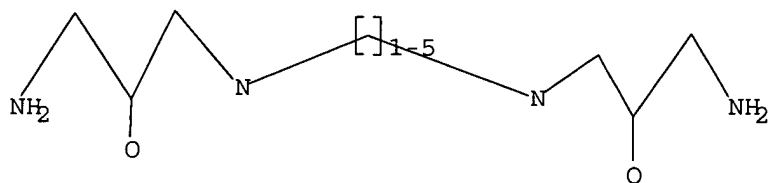
chain nodes :
 1 2 3 4 5 6 11 12 13 14 15 16 17 18 19
 chain bonds :
 1-2 1-11 1-12 2-3 3-4 4-5 5-6 11-17 12-13 13-14 14-15 15-16 15-18
 18-19
 exact/norm bonds :
 1-2 1-11 1-12 14-15 15-18 18-19
 exact bonds :
 2-3 3-4 4-5 5-6 11-17 12-13 13-14 15-16

Match level :
 1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 11:CLASS 12:CLASS 13:CLASS
 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS

L10 STRUCTURE UPLOADED

=> d query

L10 STR



Structure attributes must be viewed using STN Express query preparation.

=> s l10

SAMPLE SEARCH INITIATED 14:54:31 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 3239 TO ITERATE

30.9% PROCESSED 1000 ITERATIONS 0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 61367 TO 68193
PROJECTED ANSWERS: 0 TO 0

L11 0 SEA SSS SAM L10

=> s l10 full
FULL SEARCH INITIATED 14:54:35 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 63729 TO ITERATE

100.0% PROCESSED 63729 ITERATIONS 10 ANSWERS
SEARCH TIME: 00.00.02

L12 10 SEA SSS FUL L10

=> fil caplus		
COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	162.62	498.50
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-1.46

FILE 'CAPLUS' ENTERED AT 14:54:40 ON 25 APR 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

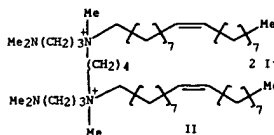
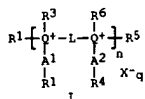
FILE COVERS 1907 - 25 Apr 2005 VOL 142 ISS 18
FILE LAST UPDATED: 24 Apr 2005 (20050424/ED)

New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l12
L13 7 L12

=> d 113 1-7 abs ibib hitstr



II

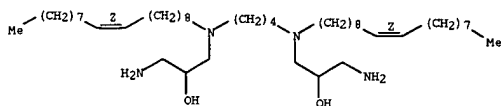
ACCESSION NUMBER:
DOCUMENT NUMBER:
TITLE:

INVENTOR(S) :
PATENT ASSIGNEE :
SOURCE :

DOCUMENT TYPE:
LANGUAGE:

FAMILY ACC. NUM.
PATENT INFORMATION

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2000027795	A1	20000518	WG 1999-UY26825	19991112
W:	AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, DE, DK, DM, EE, ES, FI, GB, GD, GE, GH, GM, GR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UY, VN, YU, ZA, ZW, AM, AZ,			
BY, GM, KE, KM, KU, RW, TD, TL				
RW:	GK, GZ, KX, LS, MW, SZ, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LJ, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, CA, GN, GW, HT, IG, HR, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UY, VN, YU, ZA, ZW, AM, AZ,			
CA 2350882	AA	20000518	CA 1999-2350882	19991112
EP 1290654	A1	20010905	EP 1999-717194	19991112
R:	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LJ, LU, NL, SE, MC, PT,			



IE, SI, LT, LV, FI, RO				
JP 2002529439	T2	20020910	JP 2000-580975	19991112
NZ 512244	A	20031219	NZ 1999-512244	19991112
AU 772847	B2	20040506	AU 2000-14776	19991112
PRIORITY APPLN. INFO.:			US 1998-108117P	P 19981112
			WO 1999-US26825	W 19991112

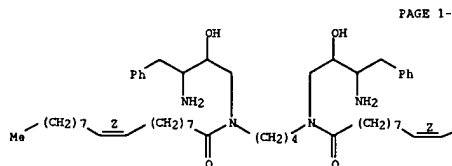
PRIORITY APPLN. INFO.:

OTHER SOURCE(S) : MARPAT 132:334312

OTHER SOURCE(S):
IT 268554-14-9P

BY 268554 34

Double bond geometry as shown.



PAGE 1-A

PAGE 1-B



IT 268539-48-6P

RN 268539-48

Double bond geometry as shown.

50-75%

ACCESSION NUMB

DOCUMENT NUMBE

AUTHOR(S) :

SOURCE:

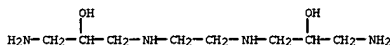
马金凤及(阿婆) 曹建琴

DOCUMENT TYPE:
LANGUAGE:

IT 58770-22-

(прера

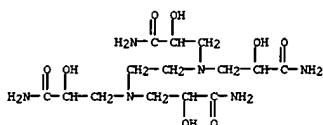
1452,



L13 ANSWER 5 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN
 AB The heat resistance of a 3:97 1,3-dioxolane-trioxane copolymer [24969-26-4] (containing 0.02% NaF and 0.2-0.4% phenolic antioxidant) is increased by adding 0.1-0.2% nitritoltrilactamide [23918-39-0] or ethylenediaminetetralactamide [23918-41-4].
 ACCESSION NUMBER: 1972:128123 CAPLUS
 DOCUMENT NUMBER: 76:128123
 TITLE: Stabilizing thermoplastic polyacetals
 INVENTOR(S): Sander, Bruno; Schuetta, Wilhelm
 PATENT ASSIGNEE(S): Badische Anilin- & Soda-Fabrik AG
 SOURCE: Ger., 3 pp.
 CODEN: GWXKAW
 DOCUMENT TYPE: Patent
 LANGUAGE: German
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
DE 1669728	A	19700924	DE 1967-B94973	19671017
DE 1669728	B	19711230		

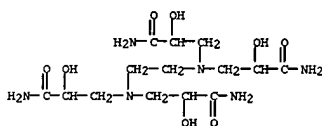
PRIORITY APPLN. INFO.: DE 1967-B94973 A 19671017
 IT 23918-41-4
 RL: PEP (Physical, engineering or chemical process); PROC (Process) (heat stabilizers, for polyoxymethylenes)
 RN 23918-41-4 CAPLUS
 CN Propanamide, 3,3',3'',3'''-(1,2-ethanediylidinitrilo)tetrakis[2-hydroxy-(9CI)] (CA INDEX NAME)



L13 ANSWER 6 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN
 AB Use of 0.1-0.2% nitritoltris(lactamide) or ethylenediaminetetrakis(lactamide) e) and 0.2-0.4% 2,2'-methylenebis(4-methyl - 6-tert-butylphenol) (I) antioxidant in degassed melts of 97:3 trioxane-1,3-dioxolane copolymer containing NaF gave 0.54-0.9% weight loss and no discoloration after heating 2 hr at 222° compared with use of I and nitritoltris(propionamide) or dicyandiamide, which gave 0.95-1.93% weight losses, brown or yellow-brown discolorations, and (or) deposits in the mold at 220-30°.
 ACCESSION NUMBER: 1970:499607 CAPLUS
 DOCUMENT NUMBER: 73:99607
 TITLE: Heat stabilization of polyacetals by addition of a β-aminolactamide
 PATENT ASSIGNEE(S): Badische Anilin- & Soda-Fabrik AG
 SOURCE: Fr., 8 pp.
 CODEN: FRXKAX
 DOCUMENT TYPE: Patent
 LANGUAGE: French
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
FR 1584931		19700102	FR	
DE 1669727			DE	
GB 1234057			GB	
US 3607831		19710000	US	

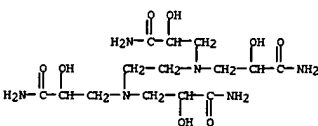
PRIORITY APPLN. INFO.: DE 19671017
 IT 23918-41-4
 RL: USES (Uses) (stabilizers, for trioxane copolymers)
 RN 23918-41-4 CAPLUS
 CN Propanamide, 3,3',3'',3'''-(1,2-ethanediylidinitrilo)tetrakis[2-hydroxy-(9CI)] (CA INDEX NAME)



L13 ANSWER 7 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN
 GI For diagram(s), see printed CA Issue.
 AB β-Carbamoyl-β-hydroxyethylamines, useful for forming complexes with Fe(III) and, to a less extent, with Ni, Co Fe(II), Cr(III), or Al, were prepared by treating glycidamide I (R = H, CH2OH, or CH2OMe) with NH3, N2H4, H2NOH, H2N(CH2)2NH2, H2NCH2CH2NH2, H2N(CH2)3NH2, H2N(CH2)6NH2, H2N(CH2CH2NH)2H, H2N(CH2CH2NH)3H, or polyethylenimine. The following HON[CH2CH(OH)CONHR]3-n were prepared (n, % yield, R, and m.p. given): 0, 86, H, 164°; 0, 76, CH2OH, oil; 1, 82, H, 193°; 1, 77, CH2OH, oil. The following [RHNCOCH(OH)CH2]2NR1N[CH2CH(OH)CONHR]2 were prepared (R1, R, % yield, and m.p. given): CH2CH2, H, 92, 201°; CHMeCH2, H, 81, oil; (CH2)3, H, 78, oil; CH2CH2, CH2 OH, 84, oil; CH2CH2, CH2OMe, 91, oil; (CH2)6, H, 91, oil; (CH2)6, CH2OH, 86, oil. Also prepared were [H2NCOCH(OH)CH2]2NN[CH2CH(OH)CONH2]2 (94%, oil); HON-[CH2CH(OH)CONH2]2 (80%, oil); HOCH2CH2N[CH2CH(OH)CONH2]2 (89%, oil); [-N[CH2CH(OH)CONH2]CH2CH2-]n (polymer) (95%, m. 60°); [H2NCOCH(OH)CH2]2N[CH2CH2NCH2CH(OH)CONH]nH, where n is 2 (88%, m. 52°) and 3 (90%, m. 48°).
 ACCESSION NUMBER: 1970:22758 CAPLUS
 DOCUMENT NUMBER: 72:22758
 TITLE: β-Carbamoyl-β-hydroxyethylamines as masking agents in the textile and paper industries
 PATENT ASSIGNEE(S): Badische Anilin- & Soda-Fabrik AG
 SOURCE: Fr., 6 pp.
 CODEN: FRXKAX
 DOCUMENT TYPE: Patent
 LANGUAGE: French
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
FR 1542198		19681011	FR	
GB 1195427			GB	
US 3538158		19700000	US	

PRIORITY APPLN. INFO.: DE 19661104
 IT 23918-41-4D, Lactamide, 3,3',3'',3'''-(ethylenedinitrilo)tetrakis-, iron complexes
 RL: USES (Uses) (as masking agents for paper)
 RN 23918-41-4 CAPLUS
 CN Propanamide, 3,3',3'',3'''-(1,2-ethanediylidinitrilo)tetrakis[2-hydroxy-(9CI)] (CA INDEX NAME)



```
=> fil reg
COST IN U.S. DOLLARS          SINCE FILE      TOTAL
                                ENTRY      SESSION
FULL ESTIMATED COST          35.48      533.98

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)  SINCE FILE      TOTAL
                                                ENTRY      SESSION
CA SUBSCRIBER PRICE          -5.11      -6.57
```

FILE 'REGISTRY' ENTERED AT 14:55:46 ON 25 APR 2005
 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
 PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
 COPYRIGHT (C) 2005 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
 provided by InfoChem.

STRUCTURE FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9
 DICTIONARY FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

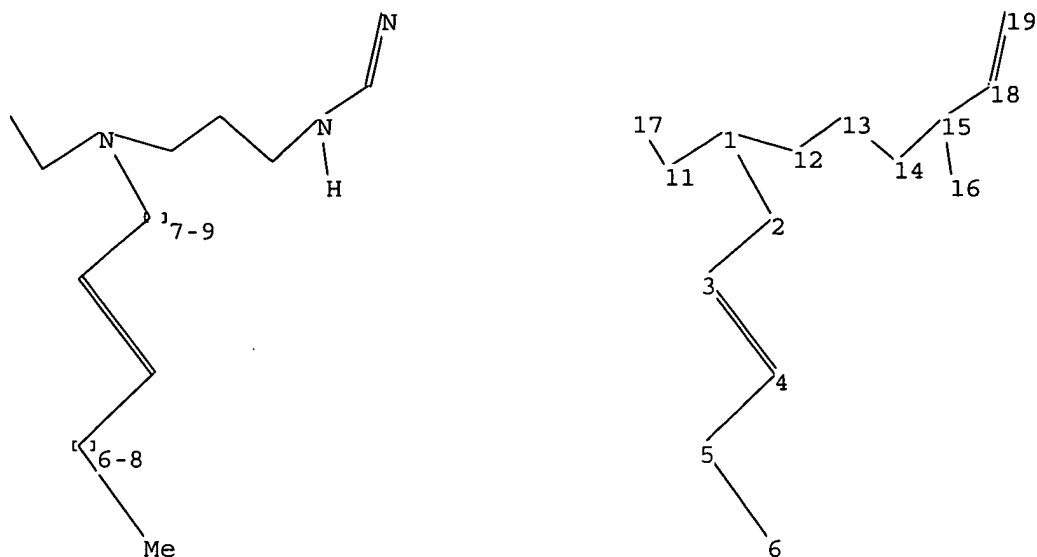
Please note that search-term pricing does apply when
 conducting SmartSELECT searches.

```
*****
*
* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added,  *
* effective March 20, 2005. A new display format, IDERL, is now    *
* available and contains the CA role and document type information. *
*
*****
```

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
 information enter HELP PROP at an arrow prompt in the file or refer
 to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

```
=>
Uploading C:\Program Files\Stnexp\Queries\09438365.str
```

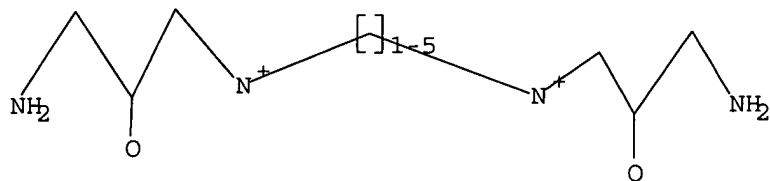
chain nodes :
 1 2 3 4 5 6 11 12 13 14 15 16 17 18 19
 chain bonds :
 1-2 1-11 1-12 2-3 3-4 4-5 5-6 11-17 12-13 13-14 14-15 15-16 15-18
 18-19
 exact/norm bonds :
 1-2 1-11 1-12 14-15 15-18 18-19
 exact bonds :
 2-3 3-4 4-5 5-6 11-17 12-13 13-14 15-16

Match level :
 1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 11:CLASS 12:CLASS 13:CLASS
 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS

L14 STRUCTURE UPLOADED

=> d query

L14 STR



Structure attributes must be viewed using STN Express query preparation.

=> s l14 full
 FULL SEARCH INITIATED 14:57:25 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 63729 TO ITERATE

100.0% PROCESSED 63729 ITERATIONS
SEARCH TIME: 00.00.01

2 ANSWERS

L15 2 SEA SSS FUL L14

=> fil caplus

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	162.19	696.17
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-6.57

FILE 'CAPLUS' ENTERED AT 14:57:29 ON 25 APR 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 25 Apr 2005 VOL 142 ISS 18
FILE LAST UPDATED: 24 Apr 2005 (20050424/ED)

New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l15

L16 0 L15

=> fil reg

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	0.45	696.62
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-6.57

FILE 'REGISTRY' ENTERED AT 14:57:40 ON 25 APR 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2005 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file

provided by InfoChem.

STRUCTURE FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9
DICTIONARY FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

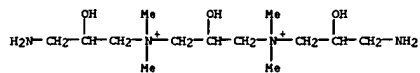
```
*****
*
* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added,   *
* effective March 20, 2005. A new display format, IDERL, is now      *
* available and contains the CA role and document type information. *
*
*****
```

Crossover limits have been increased. See HELP CROSSOVER for details.

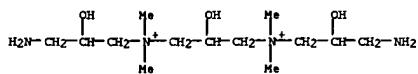
Experimental and calculated property data are now available. For more
information enter HELP PROP at an arrow prompt in the file or refer
to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=> d 115 1-2

L15 ANSWER 1 OF 2 REGISTRY COPYRIGHT 2005 ACS on STN
 RN 741631-75-4 REGISTRY
 ED Entered STN: 08 Sep 2004
 CN 1,3-Propanediaminium, N,N'-bis(3-amino-2-hydroxypropyl)-2-hydroxy-
 N,N,N',N'-tetramethyl- (9CI) (CA INDEX NAME)
 FS 3D CONCORD
 MF C13 H34 N4 O3
 CI COM
 SR CA



L15 ANSWER 2 OF 2 REGISTRY COPYRIGHT 2005 ACS on STN
 RN 93951-13-4 REGISTRY
 ED Entered STN: 31 Aug 1985
 CN 1,3-Propanediaminium, N,N'-bis(3-amino-2-hydroxypropyl)-2-hydroxy-
 N,N,N',N'-tetramethyl-, dichloride, dihydrochloride (9CI) (CA INDEX NAME)
 MF C13 H34 N4 O3 . 2 Cl H . 2 Cl
 SR European Union (EU)
 LC STN Files: CHEMLIST
 Other Sources: EINECS**
 (**Enter CHEMLIST File for up-to-date regulatory information)
 CRN (741631-75-4)



● 2 Cl⁻

● 2 HCl

=> FIL STNGUIDE
COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
4.54	701.16

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
0.00	-6.57

CA SUBSCRIBER PRICE

FILE 'STNGUIDE' ENTERED AT 14:58:55 ON 25 APR 2005
USE IS SUBJECT TO THE TERMS OF YOUR CUSTOMER AGREEMENT
COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY, JAPAN SCIENCE
AND TECHNOLOGY CORPORATION, AND FACHINFORMATIONSZENTRUM KARLSRUHE

FILE CONTAINS CURRENT INFORMATION.
LAST RELOADED: Apr 22, 2005 (20050422/UP).

=> fil reg
COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
0.06	701.22

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
0.00	-6.57

CA SUBSCRIBER PRICE

FILE 'REGISTRY' ENTERED AT 14:59:06 ON 25 APR 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2005 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
provided by InfoChem.

STRUCTURE FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9
DICTIONARY FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

*
* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added, *
* effective March 20, 2005. A new display format, IDERL, is now *
* available and contains the CA role and document type information. *
*

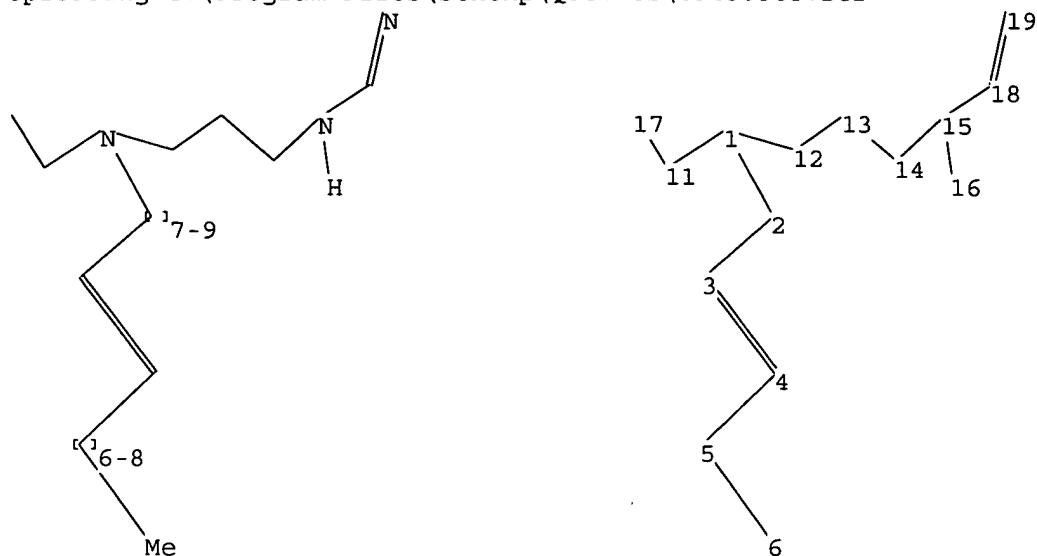
Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
information enter HELP PROP at an arrow prompt in the file or refer
to the file summary sheet on the web at:

<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>

Uploading C:\Program Files\Stnexp\Queries\09438365.str



chain nodes :

1 2 3 4 5 6 11 12 13 14 15 16 17 18 19

chain bonds :

1-2 1-11 1-12 2-3 3-4 4-5 5-6 11-17 12-13 13-14 14-15 15-16 15-18 18-19

exact/norm bonds :

1-2 1-11 1-12 14-15 15-18 18-19

exact bonds :

2-3 3-4 4-5 5-6 11-17 12-13 13-14 15-16

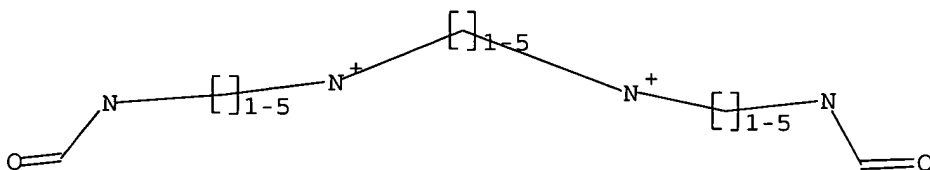
Match level :

1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 11:CLASS 12:CLASS 13:CLASS
14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS

L17 STRUCTURE UPLOADED

=> d query

L17 STR



Structure attributes must be viewed using STN Express query preparation.

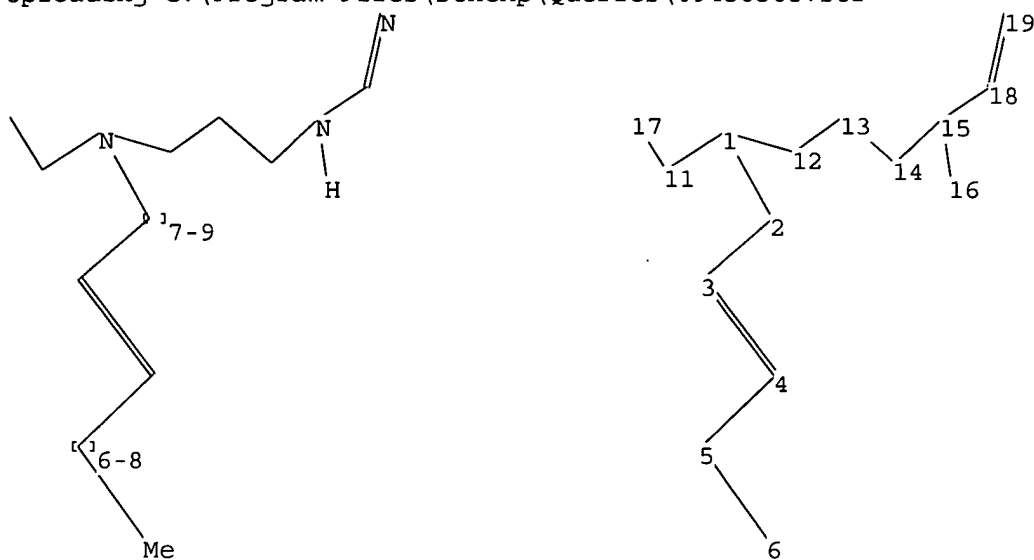
=> s l17
 SAMPLE SEARCH INITIATED 15:01:38 FILE 'REGISTRY'
 SAMPLE SCREEN SEARCH COMPLETED - 62612 TO ITERATE

1.6% PROCESSED 1000 ITERATIONS 0 ANSWERS
 INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
 SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**
 BATCH **INCOMPLETE**
 PROJECTED ITERATIONS: EXCEEDS 1000000
 PROJECTED ANSWERS: EXCEEDS 0

L18 0 SEA SSS SAM L17

=>
 Uploading C:\Program Files\Stnexp\Queries\09438365.str



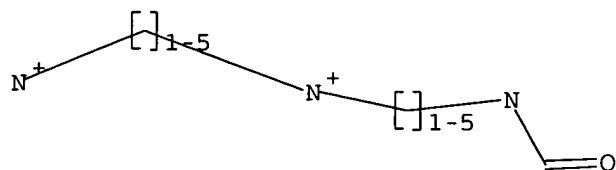
chain nodes :
 1 2 3 4 5 6 11 12 13 14 15 16 17 18 19
 chain bonds :
 1-2 1-11 1-12 2-3 3-4 4-5 5-6 11-17 12-13 13-14 14-15 15-16 15-18
 18-19
 exact/norm bonds :
 1-2 1-11 1-12 14-15 15-18 18-19
 exact bonds :
 2-3 3-4 4-5 5-6 11-17 12-13 13-14 15-16

Match level :
 1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 11:CLASS 12:CLASS 13:CLASS
 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS

L19 STRUCTURE UPLOADED

=> d query

L19 STR



Structure attributes must be viewed using STN Express query preparation.

=> s 119

SAMPLE SEARCH INITIATED 15:02:18 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 86223 TO ITERATE

1.2% PROCESSED 1000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

1 ANSWERS

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**
 BATCH **INCOMPLETE**

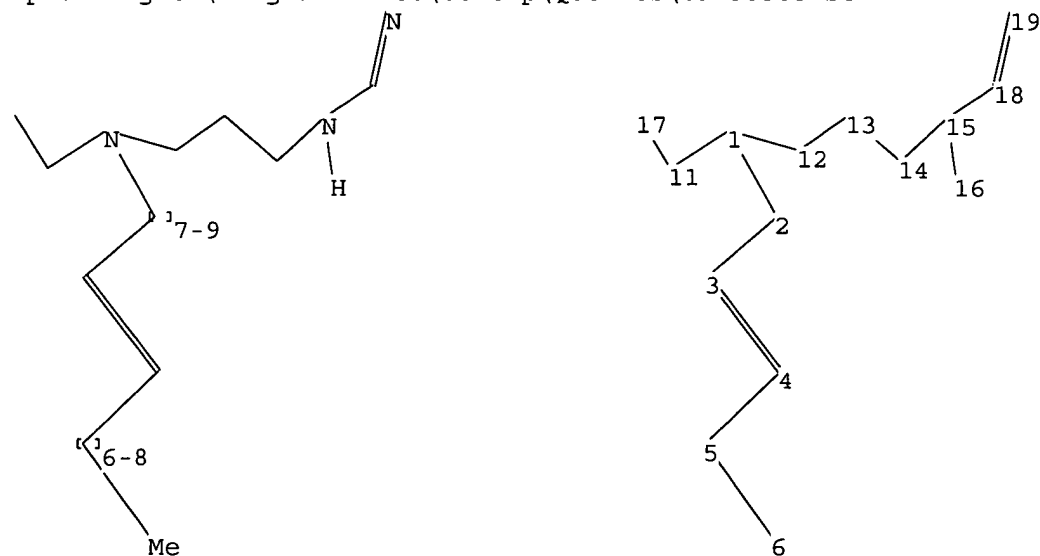
PROJECTED ITERATIONS: EXCEEDS 1000000

PROJECTED ANSWERS: EXCEEDS 1167

L20 1 SEA SSS SAM L19

=>

Uploading C:\Program Files\Stnexp\Queries\09438365.str



chain nodes :

1 2 3 4 5 6 11 12 13 14 15 16 17 18 19

chain bonds :

1-2 1-11 1-12 2-3 3-4 4-5 5-6 11-17 12-13 13-14 14-15 15-16 15-18
18-19

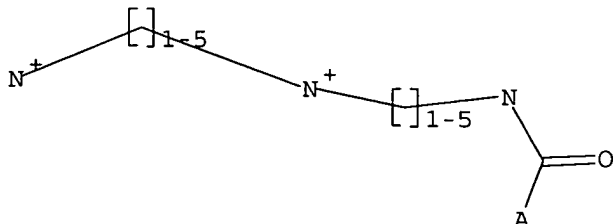
exact/norm bonds :
1-2 1-11 1-12 14-15 15-18 18-19
exact bonds :
2-3 3-4 4-5 5-6 11-17 12-13 13-14 15-16

Match level :
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 11:CLASS 12:CLASS 13:CLASS
14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS

L21 STRUCTURE UPLOADED

=> d query

L21 STR



Structure attributes must be viewed using STN Express query preparation.

=> s l21

SAMPLE SEARCH INITIATED 15:03:17 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 86218 TO ITERATE

1.2% PROCESSED 1000 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

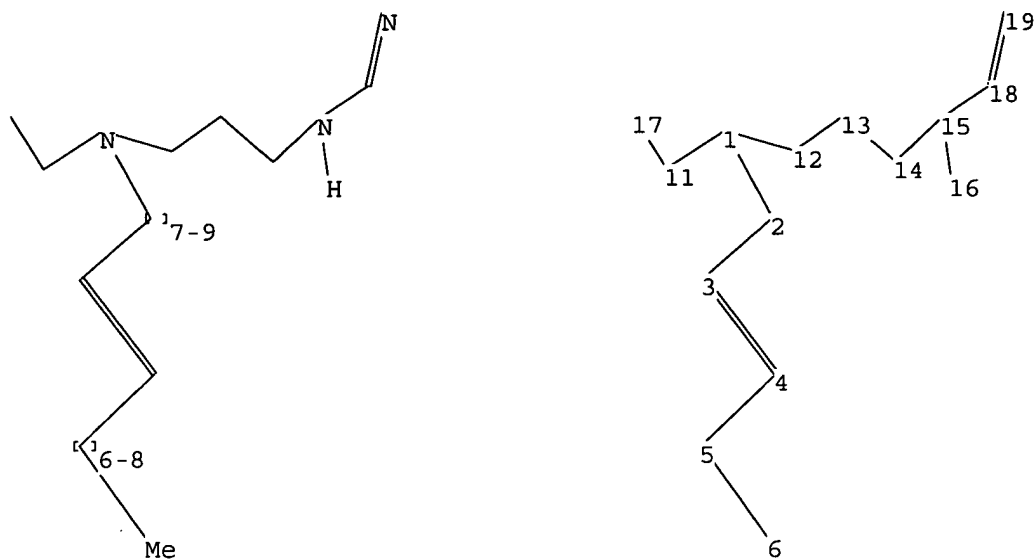
0 ANSWERS

FULL FILE PROJECTIONS: ONLINE **INCOMPLETE**
BATCH **INCOMPLETE**
PROJECTED ITERATIONS: EXCEEDS 1000000
PROJECTED ANSWERS: EXCEEDS 0

L22 0 SEA SSS SAM L21

=>

Uploading C:\Program Files\Stnexp\Queries\09438365.str



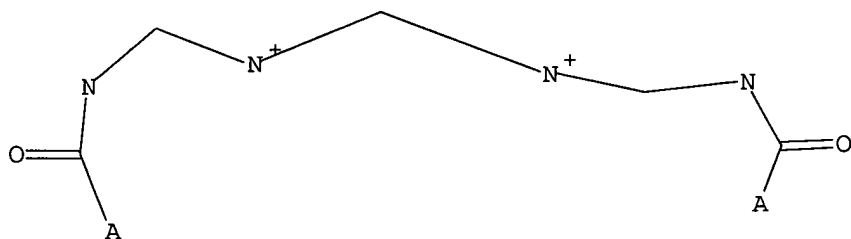
chain nodes :
 1 2 3 4 5 6 11 12 13 14 15 16 17 18 19
 chain bonds :
 1-2 1-11 1-12 2-3 3-4 4-5 5-6 11-17 12-13 13-14 14-15 15-16 15-18
 18-19
 exact/norm bonds :
 1-2 1-11 1-12 14-15 15-18 18-19
 exact bonds :
 2-3 3-4 4-5 5-6 11-17 12-13 13-14 15-16

Match level :
 1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 11:CLASS 12:CLASS 13:CLASS
 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS

L23 STRUCTURE UPLOADED

=> d query

L23 STR



Structure attributes must be viewed using STN Express query preparation.

=> s 123
 SAMPLE SEARCH INITIATED 15:04:26 FILE 'REGISTRY'
 SAMPLE SCREEN SEARCH COMPLETED - 68 TO ITERATE

100.0% PROCESSED 68 ITERATIONS 0 ANSWERS
 SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
 BATCH **COMPLETE**
 PROJECTED ITERATIONS: 866 TO 1854
 PROJECTED ANSWERS: 0 TO 0

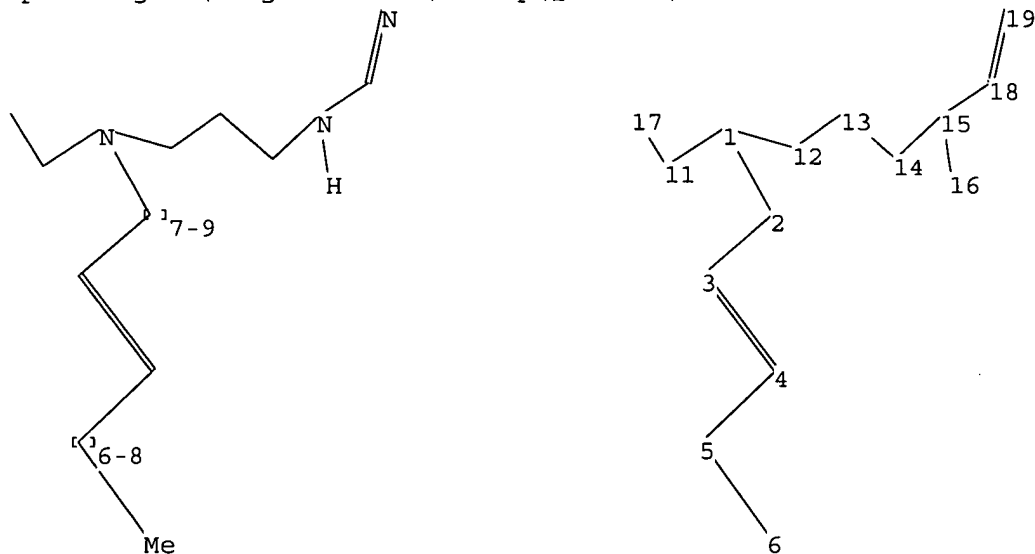
L24 0 SEA SSS SAM L23

=> s 123 full
 FULL SEARCH INITIATED 15:04:30 FILE 'REGISTRY'
 FULL SCREEN SEARCH COMPLETED - 1654 TO ITERATE

100.0% PROCESSED 1654 ITERATIONS 0 ANSWERS
 SEARCH TIME: 00.00.01

L25 0 SEA SSS FUL L23

=>
 Uploading C:\Program Files\Stnexp\Queries\09438365.str



chain nodes :
 1 2 3 4 5 6 11 12 13 14 15 16 17 18 19
 chain bonds :
 1-2 1-11 1-12 2-3 3-4 4-5 5-6 11-17 12-13 13-14 14-15 15-16 15-18
 18-19
 exact/norm bonds :
 1-2 1-11 1-12 14-15 15-18 18-19
 exact bonds :
 2-3 3-4 4-5 5-6 11-17 12-13 13-14 15-16

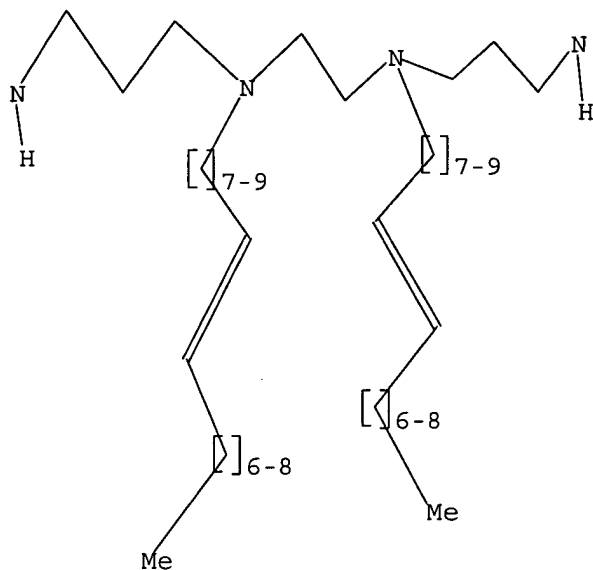
Match level :

1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 11:CLASS 12:CLASS 13:CLASS
14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS

L26 STRUCTURE UPLOADED

=> d query

L26 STR



Structure attributes must be viewed using STN Express query preparation.

=> s l26

SAMPLE SEARCH INITIATED 15:08:28 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 17 TO ITERATE

100.0% PROCESSED 17 ITERATIONS
SEARCH TIME: 00.00.01

1 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 93 TO 587
PROJECTED ANSWERS: 1 TO 80

L27 1 SEA SSS SAM L26

=> s l26 full

FULL SEARCH INITIATED 15:08:34 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 257 TO ITERATE

100.0% PROCESSED 257 ITERATIONS
SEARCH TIME: 00.00.01

4 ANSWERS

L28 4 SEA SSS FUL L26

=> fil caplus

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

328.68

1029.90

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

ENTRY

SESSION

CA SUBSCRIBER PRICE

0.00

-6.57

FILE 'CAPLUS' ENTERED AT 15:08:37 ON 25 APR 2005

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 25 Apr 2005 VOL 142 ISS 18

FILE LAST UPDATED: 24 Apr 2005 (20050424/ED)

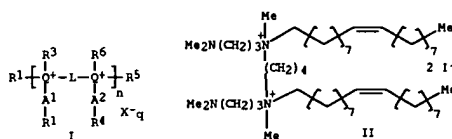
New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l28

L29 2 L28

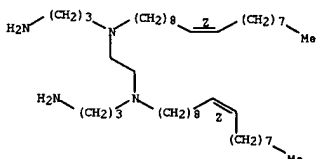
=> d l29 1-2 abs ibib hitstr



AB Synthesis and activity of transfection reagents (I) [Q = N, O, S; L = (un)substituted alkyl, ether, polyether, amide, polyamide, ester, sulfide, urea, thiourea, guanidyl, carbamoyl, carbonate, phosphate, sulfate, sulfoxide, imine, carbonyl, secondary amine; R1-R6 independently = (un)substituted alkyl, alkenyl, aryl, ether; A1, A2 independently = CH2O, CH2S, CH2NH, CO, C=NH, CS, alkyl; X = physiol. acceptable anion; n = 1-1000; q = number of pos. charge divided by valence of anion], cationic lipids capable of facilitating transport of biol. active agents or substances into cells, are disclosed. Thus, I [R1,R4 = oleyl; R2,R5 = Me2N(CH2)3; R3,R6 = Me; A1,A2 = CH2; L = (CH2)4; X = I] (II) is prepared by reaction of bis-1,4-oleyl-1,4-butanediamine with acrylonitrile followed by reduction of nitrile to amine and quaternization of amine with Me iodide.

II shows an activity of 37.8 ng/µg/cm2 in DNA delivery. Formulations containing I are given.
ACCESSION NUMBER: 2000:335366 CAPLUS
DOCUMENT NUMBER: 132:334312
TITLE: synthesis and activity of transfection reagents for transport of biol. active agents or substances into cells
INVENTOR(S): Chu, Yongliang; Masoud, Malek; Gebeyehu, Gulilat
PATENT ASSIGNEE(S): Life Technologies, Inc., USA
SOURCE: PCT Int. Appl., 130 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

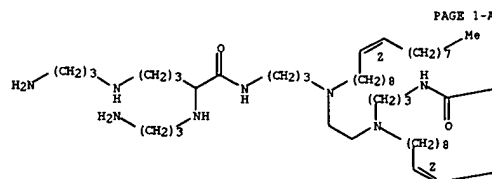
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2000027795	A1	20000518	WO 1999-US26825	19991112
W: AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, DE, DK, DM, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, A2, BY, KG, KZ, MD, RU, TJ, TM				
RW: GH, GM, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
CA 2350882	AA	20000518	CA 1999-2350882	19991112



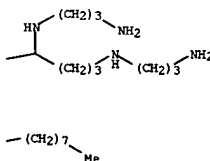
REFERENCE COUNT: 13 THERE ARE 13 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

OTHER SOURCE(S): MARPAT 132:334312
IT 268539-54-4P 268539-55-5P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(synthesis and activity of transfection reagents for transport of biol. active agents or substances into cells)
RN 268539-54-4 CAPLUS
CN Pentanamide, N,N'-[1,2-ethanediylbis[(9Z)-9-octadecenylimino]-3,1-propanediyl]]bis(2,5-bis[(3-aminopropyl)amino]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



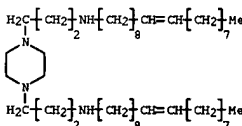
PAGE 1-A



PAGE 1-B

RN 268539-55-5 CAPLUS
CN 1,3-Propanediamine, N,N'-1,2-ethanediylbis[N-(9Z)-9-octadecenyl]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



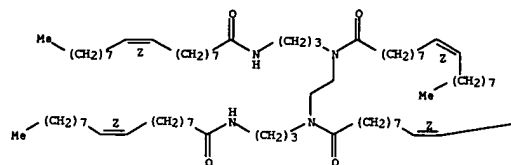
I

AB Transfection of a cell is accomplished using with a polynucleotide mixed with one or more amphipathic compds. and a DNA-binding protein, especially a histone such as histones H1, H2A, or H2B. The DNA-binding protein may be fused to a nuclear localization signal peptide. Exemplary and preferred amphipathic compds. are cationic amphipathic compds. I was synthesized in 70% yield by reacting 1,4-bis(3-aminopropyl)piperazine with oleoyl chloride and reducing the intermediate with LiAlH4 in THF. Histone H1 was found to increase the transfection efficiency of I 16.1-fold. I/H1 reagent has a greater transfection efficiency and less cellular toxicity than LipofectAmine, which is useful in gene therapy.

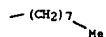
ACCESSION NUMBER: 1998:263206 CAPLUS
DOCUMENT NUMBER: 128:266964
TITLE: Process of transfecting a cell with a polynucleotide mixed with an amphipathic compound and a DNA-binding protein
INVENTOR(S): Wolff, Jon A.; Fritz, Jeffery; Budker, Vladimir; Hagstrom, James
PATENT ASSIGNEE(S): Mirus Corporation, USA
SOURCE: U.S., 16 pp.
CODEN: USXXAM
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 5744335	A	19980428	US 1995-530598	19950919
US 6180784	B1	20010130	US 1998-20566	19980117
PRIORITY APPL. INFO.:			US 1995-530598	A3 19950919
OTHER SOURCE(S): MARPAT 128:266964				
IT 205596-14-1P				
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)				
(amphipathic compds. for transfecting cells and their syntheses)				
RN 205596-14-1 CAPLUS				
CN 9-Octadecenamide, N,N'-1,2-ethanediylbis[N-[3-[(1-oxo-9-octadecenyl)amino]propyl]-, (all-Z)- (9CI) (CA INDEX NAME)]				

Double bond geometry as shown.

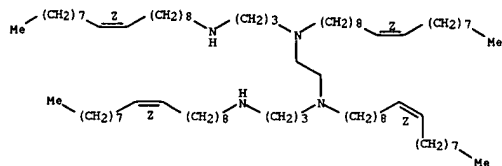


PAGE 1-B



IT 205596-06-1P
RL: BUU (Biological use, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(cell transfection with polynucleotide mixed with amphipathic compound and DNA-binding protein)
RN 205596-06-1 CAPLUS
CN 1,3-Propanediamine, N,N''-1,2-ethanediylbis[N,N'-di-9-octadecenyl-, (all-2)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



REFERENCE COUNT: 25 THERE ARE 25 CITED REFERENCES AVAILABLE FOR THIS

```
=>  fil reg
COST IN U.S. DOLLARS          SINCE FILE      TOTAL
                                ENTRY      SESSION
FULL ESTIMATED COST          11.23      1041.13

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)  SINCE FILE      TOTAL
                                                ENTRY      SESSION
CA SUBSCRIBER PRICE          -1.46      -8.03
```

FILE 'REGISTRY' ENTERED AT 15:10:19 ON 25 APR 2005
 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
 PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
 COPYRIGHT (C) 2005 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
 provided by InfoChem.

STRUCTURE FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9
 DICTIONARY FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

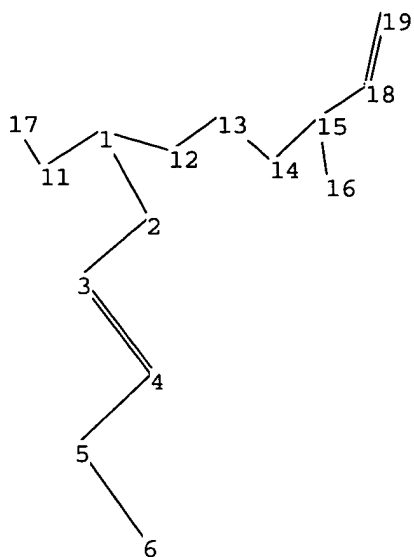
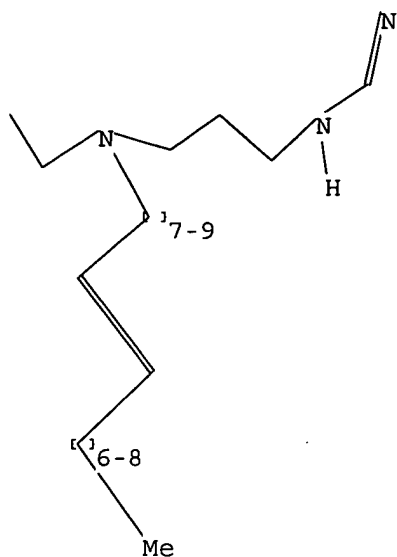
Please note that search-term pricing does apply when
 conducting SmartSELECT searches.

```
*****
*
* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added,  *
* effective March 20, 2005. A new display format, IDERL, is now    *
* available and contains the CA role and document type information. *
*
*****
```

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
 information enter HELP PROP at an arrow prompt in the file or refer
 to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

```
=>
Uploading C:\Program Files\Stnexp\Queries\09438365.str
```

```

chain nodes :
1  2  3  4  5  6  11 12 13 14 15 16 17 18 19
chain bonds :
1-2 1-11 1-12 2-3 3-4 4-5 5-6 11-17 12-13 13-14 14-15 15-16 15-18
18-19
exact/norm bonds :
1-2 1-11 1-12 14-15 15-18 18-19
exact bonds :
2-3 3-4 4-5 5-6 11-17 12-13 13-14 15-16

```

```

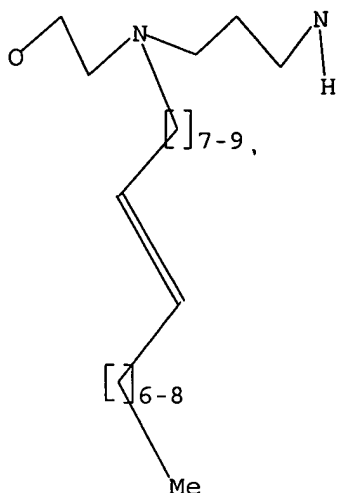
Match level :
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 11:CLASS 12:CLASS 13:CLASS
14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS

```

L30 STRUCTURE UPLOADED

=> d query

L30 STR



Structure attributes must be viewed using STN Express query preparation.

=> s l30

SAMPLE SEARCH INITIATED 15:22:13 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 62 TO ITERATE

100.0% PROCESSED 62 ITERATIONS 1 ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 768 TO 1712
PROJECTED ANSWERS: 1 TO 80

L31 1 SEA SSS SAM L30

=> s l30 full

FULL SEARCH INITIATED 15:22:17 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 1299 TO ITERATE

100.0% PROCESSED 1299 ITERATIONS 11 ANSWERS
SEARCH TIME: 00.00.01

L32 11 SEA SSS FUL L30

=> fil caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	169.50	1210.63
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-8.03

FILE 'CAPLUS' ENTERED AT 15:22:20 ON 25 APR 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 25 Apr 2005 VOL 142 ISS 18
FILE LAST UPDATED: 24 Apr 2005 (20050424/ED)

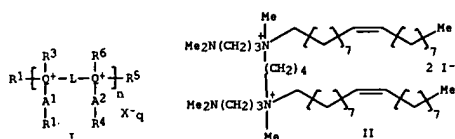
New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 132

L33 4 L32

=> d 133 1-4 abs ibib hitstr



AB Synthesis and activity of transfection reagents (I) [Q = N, O, S; L = (un)substituted alkyl, ether, polyether, amide, polyamide, ester, sulfide, urea, thiourea, guanidyl, carbamoyl, carbonate, phosphate, sulfate, sulfoxide, imine, carbonyl, secondary amine; R1-R6 independently = (un)substituted alkyl, alkenyl, aryl, ether; A1, A2 independently = CH2O, CH2S, CH2NH, CO, C=NH, CS, alkyl; X = physiol. acceptable anion; n = 1-1000; q = number of pos. charge divided by valence of anion], cationic lipids capable of facilitating transport of biol. active agents or substances into cells, are disclosed. Thus, I [R1, R4 = oleyl; R2, R5 = Me2N(CH2)3; R3, R6 = Me; A1, A2 = CH2; L = (CH2)4; X = I] (II) is prepared by reaction of bis-1,4-oleyl-1,4-butandiamine with acrylonitrile followed by reduction of nitrile to amine and quaternization of amine with Me iodide.

II shows an activity of 37.8 ng/μg/cm2 in DNA delivery. Formulations containing I are given.

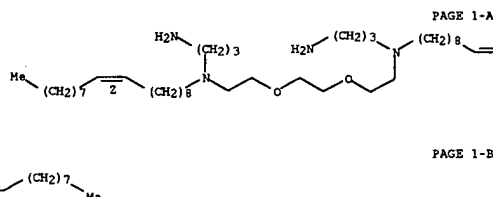
ACCESSION NUMBER: 2000:335366 CAPLUS
DOCUMENT NUMBER: 132:334312
TITLE: synthesis and activity of transfection reagents for transport of biol. active agents or substances into cells
INVENTOR(S): Chu, Yongliang; Masoud, Malek; Gebeyehu, Gulilat
PATENT ASSIGNEE(S): Life Technologies, Inc., USA
SOURCE: PCT Int. Appl., 130 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2000027795	A1	20000518	WO 1999-US26825	19991112
W:	AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, DE, DK, DM, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM			
RW:	GH, GH, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CH, GA, GN, GW, ML, MR, NE, SN, TD, TG			
CA 2350882	AA	20000518	CA 1999-2350882	19991112

L33 ANSWER 1 OF 4 CAPLUS COPYRIGHT 2005 ACS on STN (Continued)
EP 1129064 A1 20010905 EP 1999-971794 19991112
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO
JP 2002529439 T2 20020910 JP 2000-580975 19991112
NZ 512244 A 20031219 NZ 1999-512244 19991112
AU 772847 B2 20040506 AU 2000-14776 19991112
PRIORITY APPL. INFO.: US 1998-108117P P 19981112
WO 1999-US26825 W 19991112

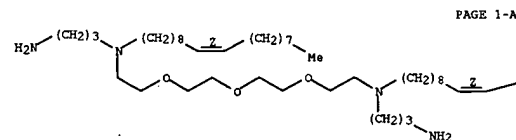
OTHER SOURCE(S): MARPAT 132:334312
IT 268539-56-6P 268539-57-7P 268539-58-8P
268539-60-2P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(synthesis and activity of transfection reagents for transport of biol. active agents or substances into cells)
RN 268539-56-6 CAPLUS
CN 7,10-Dioxo-4,13-diazahexadecane-1,16-diamine, 4,13-di-(9Z)-9-octadecenyl- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

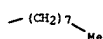


RN 268539-57-7 CAPLUS
CN 7,10,13-Trioxo-4,16-diazanonadecane-1,19-diamine, 4,16-di-(9Z)-9-octadecenyl- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



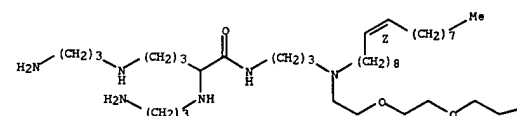
PAGE 1-B



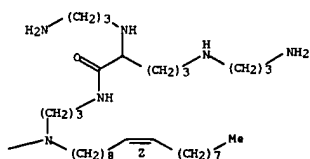
RN 268539-58-8 CAPLUS
CN Pentanamide, N,N'-[4,13-di-(9Z)-9-octadecenyl-7,10-dioxo-4,13-diazahexadecane-1,16-diyl]bis[2,5-bis[(3-aminopropyl)amino]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

PAGE 1-A



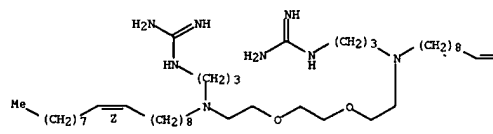
PAGE 1-B



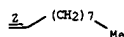
RN 268539-60-2 CAPLUS
CN 9,12-Dioxo-2,6,15,19-tetraazaoicosanediimidamide, 6,15-di-(9Z)-9-octadecenyl- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

PAGE 1-A



PAGE 1-B



REFERENCE COUNT: 13 THERE ARE 13 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L33 ANSWER 2 OF 4 CAPLUS COPYRIGHT 2005 ACS on STN
 AB ABCD[N]([CH2]nOXY) ([CH2]oOXY) [A = H, NR1R2, NR1(CH2)pNR3R4, (C:R)NH2, pyridinyl; B, D = bond, C1-6 alkylene, iminoalkylene; C = piperidinediyl, piperazinediyl; W, X = bond, CO; Y, Z = (unsatd.) hydrocarbyl; R1-R5 = H, alkyl; n = 0-2; o, p = 2-6; with provisos], were prepared Thus, 2-[(2-hydroxyethyl)-[1-(3-diethylaminopropyl)piperidin-4-ylmethyl]amino]ethanol reacted with tetradecanoyl chloride to give 314 tetradecanoic acid 2-[(2-tetradecanoyloxyethyl)-[1-(3-diethylaminopropyl)piperidin-4-ylmethyl]amino]ethyl ester. The latter was 3.2 times as effective as DOTAP for transfection of DNA into HeLa cells.

ACCESSION NUMBER: 1998:180856 CAPLUS
 DOCUMENT NUMBER: 128:230250
 TITLE: Preparation of (heterocyclic) amino alcohol derivatives as agents for introducing bioactive molecules into cells.

INVENTOR(S): Friebe, Walter-Gunar; Dimoudis, Nikolaos; Michaelis, Uwe; Knipp, Bernhard
 PATENT ASSIGNEE(S): Boehringer Mannheim G.m.b.H., Germany; Friebe, Walter-Gunar; Dimoudis, Nikolaos; Michaelis, Uwe; Knipp, Bernhard
 SOURCE: PCT Int. Appl., 71 pp.
 CODEN: PIXX02
 DOCUMENT TYPE: Patent
 LANGUAGE: German
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9811082	A1	19980319	WO 1997-EP4944	19970910
W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GR, HU, ID, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
RW: GH, KE, LS, MW, SD, SZ, UG, ZW, AT, BE, CH, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, ML, MR, NE, SN, TD, TG				
DE 19637043	A1	19980319	DE 1996-19637043	19960912
CA 2265765	AA	19980319	CA 1997-2265765	19970910
AU 9743841	A1	19980402	AU 1997-43841	19970910
AU 741636	B2	20011206		
EP 927174	A1	19990707	EP 1997-942012	19970910
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, PT, IE, FI				
CN 1237161	A	19991201	CN 1997-199651	19970910
CN 1087293	B	20020710		
BR 9712818	A	19991221	BR 1997-12818	19970910
JP 2001504806	T2	20010410	JP 1998-513249	19970910
ZA 9708170	A	19990311	ZA 1997-8170	19970911
KR 2000036063	A	20000626	KR 1999-702068	19990311
US 2003236266	A1	20031225	US 2002-59207	20020131
PRIORITY APPLN. INFO.:			DE 1996-19637043	A 19960912
			WO 1997-EP4944	W 19970910
			US 1999-147818	B1 19990512

OTHER SOURCE(S): MARPAT 128:230250
 IT 204639-32-7P
 RI: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses).
 (preparation of (heterocyclic) amino alc. derivs. as agents for introducing

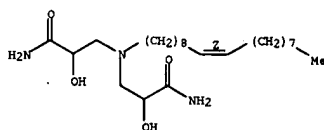
L33 ANSWER 3 OF 4 CAPLUS COPYRIGHT 2005 ACS on STN
 AB The title compds. are useful as emulsifying agents and as fat-liquoring or softening agents for leather, paper, and textiles. Inter-medates RNHCH2CH(OH)CONH2 (I) and RN[CH2CH(OH)-CONH2]2 (II) were prepared by reacting 1 or 2 moles glycidamide with 1 mole of amine RNH2 in 25 aqueous, aqueous methanolic, or methanolic solution at 50-60° during 1-2 hr. Quaternization of I and II was carried out by treating with an appropriate halide in H2O, aqueous methanolic or methanolic solution at 30-50° for 2-3 hr then at 50° for 2-3 hr. Aqueous solns. containing 0.3 g/l. [C18H39-NMe[CH2CH(OH)CONH2]2]+MeSO4- were used to treat cotton and polycaprolactam fabrics giving material with a soft pleasant hand and hydrophilic in character.

ACCESSION NUMBER: 1971:87345 CAPLUS
 DOCUMENT NUMBER: 74:87345
 TITLE: (β-Carbamyl-β-hydroxyethyl)alkylammonium salts
 PATENT ASSIGNEE(S): Badische Anilin- & Soda-Fabrik AG
 SOURCE: Brit., 12 pp.
 CODEN: BRXKAA
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
GB 1211040		19701104	GB	
FR 1592740			FR	
US 3632623		19720000	US	
PRIORITY APPLN. INFO.:			DE	19671121

IT 32671-73-1P 32671-74-2P
 RL: SPN (Synthetic preparation); PREP (Preparation) (preparation of)
 RN 32671-73-1 CAPLUS
 CN Lactamide, 3,3'-(9-octadecenylimino)bis-, (Z)- (8CI) (CA INDEX NAME)

Double bond geometry as shown.

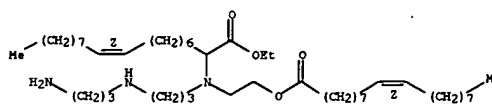


RN 32671-74-2 CAPLUS
 CN Ammonium, bis(2-carbamyl-2-hydroxyethyl)methyl-9-octadecenyl-, methyl sulfate, (Z)- (8CI) (CA INDEX NAME)

CM 1
 CRN 48077-37-8
 CMF C25 H50 N3 O4

L33 ANSWER 2 OF 4 CAPLUS COPYRIGHT 2005 ACS on STN (Continued)
 bioactive molcs. into cells)
 RN 204639-32-7 CAPLUS
 CN 9-Octadecenoic acid, 2-[[[3-[(3-aminopropyl)amino]propyl][2-[[[1-oxo-9-octadecenyl)oxy]ethyl]amino]-, ethyl ester, monohydrochloride, (Z,Z)- (9CI) (CA INDEX NAME)

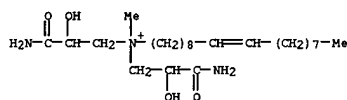
Double bond geometry as shown.



● HCl

REFERENCE COUNT: 19 THERE ARE 19 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L33 ANSWER 3 OF 4 CAPLUS COPYRIGHT 2005 ACS on STN (Continued)



CM 2

CRN 21228-90-0
 CMF C H3 O4 S

Me-O-SO3-

AB Bituminous emulsions for road building are described which break apart quickly. These neutral or slightly acid cationic aqueous bituminous emulsions

are prepared with N-alkylpropylenediamine salts as emulsifiers. They contain 20-80% bitumen, 0.1-2% emulsifier C8-22H7-45N*H-[(CH₂CH₂O)₁₁-(CH₂)₃NHCH₂], while the rest is water. The pH is 6.5-7.0. The diamine is prepared by letting a β-cyano-ethylalkylamine react with 1-4 moles ethylene oxide to yield a polyoxyethylenated β-cyanoethylalkylamine. The latter is catalytically hydrogenated to give the diamine. Thus, 35% water and 0.1% N-oleyl-N-(hydroxyethyl)-1,3-propylenediamine yielded an aqueous solution of an emulsifier which subsequently was mixed with 65%

asphalt

to result in an emulsion which broke quickly.

ACCESSION NUMBER: 1969:423534 CAPLUS

DOCUMENT NUMBER: 71:23534

TITLE: Aqueous bituminous emulsions

INVENTOR(S): Reck, Richard A.; Stalioraitis, Joseph S.; Dybalski, Jack N.

PATENT ASSIGNEE(S): Armour and Co.

SOURCE: Ger., 4 pp.

CODEN: GWXKAW

DOCUMENT TYPE: Patent

LANGUAGE: German

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
DE 1292572		19690410	US	19640511

PRIORITY APPLN. INFO.:

IT 26351-33-7

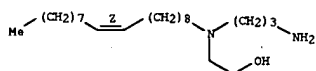
RL: USES (Uses)

(pavement containing bitumens and, cationic)

RN 26351-33-7 CAPLUS

CN Ethanol, 2-[(3-aminopropyl)-9-octadecenylamino]- (8CI) (CA INDEX NAME)

Double bond geometry as shown.



```
=> fil reg
COST IN U.S. DOLLARS          SINCE FILE      TOTAL
                                ENTRY      SESSION
FULL ESTIMATED COST          21.56      1232.19

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)  SINCE FILE      TOTAL
                                                ENTRY      SESSION
CA SUBSCRIBER PRICE          -2.92      -10.95
```

FILE 'REGISTRY' ENTERED AT 15:24:43 ON 25 APR 2005
 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
 PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
 COPYRIGHT (C) 2005 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file
 provided by InfoChem.

STRUCTURE FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9
 DICTIONARY FILE UPDATES: 24 APR 2005 HIGHEST RN 849094-71-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

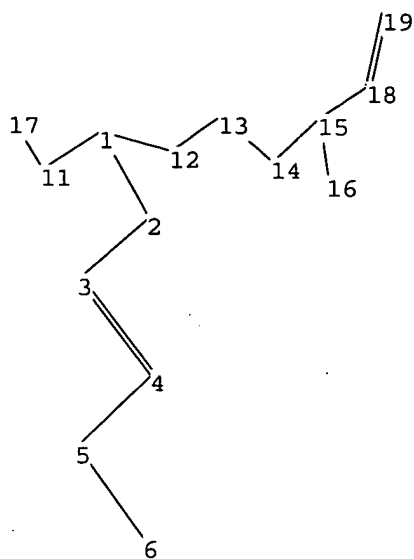
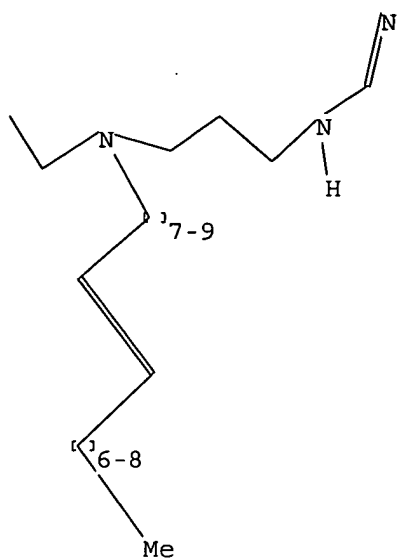
Please note that search-term pricing does apply when
 conducting SmartSELECT searches.

```
*****
*
* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added, *
* effective March 20, 2005. A new display format, IDERL, is now *
* available and contains the CA role and document type information. *
*
*****
```

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
 information enter HELP PROP at an arrow prompt in the file or refer
 to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

```
=>
Uploading C:\Program Files\Stnexp\Queries\09438365.str
```



```

chain nodes :
1 2 3 4 5 6 11 12 13 14 15 16 17 18 19
chain bonds :
1-2 1-11 1-12 2-3 3-4 4-5 5-6 11-17 12-13 13-14 14-15 15-16 15-18
18-19
exact/norm bonds :
1-2 1-11 1-12 14-15 15-18 18-19
exact bonds :
2-3 3-4 4-5 5-6 11-17 12-13 13-14 15-16

```

```

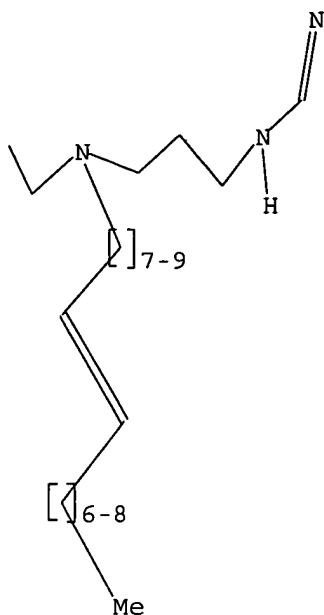
Match level :
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 11:CLASS 12:CLASS 13:CLASS
14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS

```

L34 STRUCTURE UPLOADED

=> d query

L34 STR



Structure attributes must be viewed using STN Express query preparation.

=> s 134

SAMPLE SEARCH INITIATED 15:25:29 FILE 'REGISTRY'
 SAMPLE SCREEN SEARCH COMPLETED - 135 TO ITERATE

100.0% PROCESSED 135 ITERATIONS 0 ANSWERS
 SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
 BATCH **COMPLETE**

PROJECTED ITERATIONS: 2003 TO 3397
 PROJECTED ANSWERS: 0 TO 0

L35 0 SEA SSS SAM L34

=> s 134 full

FULL SEARCH INITIATED 15:25:33 FILE 'REGISTRY'
 FULL SCREEN SEARCH COMPLETED - 2564 TO ITERATE

100.0% PROCESSED 2564 ITERATIONS 19 ANSWERS
 SEARCH TIME: 00.00.01

L36 19 SEA SSS FUL L34

=> fil caplus

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
161.33	1393.52

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION

CA SUBSCRIBER PRICE

0.00

-10.95

FILE 'CAPLUS' ENTERED AT 15:25:36 ON 25 APR 2005
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2005 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 25 Apr 2005 VOL 142 ISS 18
FILE LAST UPDATED: 24 Apr 2005 (20050424/ED)

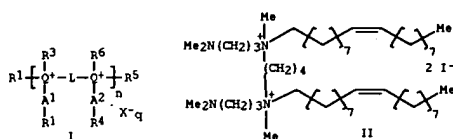
New CAS Information Use Policies, enter HELP USAGETERMS for details.

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l36

L37 3 L36

=> d l37 1-3 abs ibib hitstr



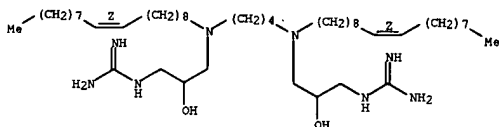
AB Synthesis and activity of transfection reagents (I) [Q = N, O, S; L = (un)substituted alkyl, ether, polyether, amide, polyamide, ester, sulfide, urea, thiourea, guanidyl, carbamoyl, carbonate, phosphate, sulfate, sulfoxide, imine, carbonyl, secondary amine; R1-R6 independently = (un)substituted alkyl, alkenyl, aryl, ether; A1, A2 independently = CH2O, CH2S, CH2NH, CO, C=NH, CS, alkyl; X = physiol. acceptable anion; a = 1-1000; q = number of pos. charge divided by valence of anion], cationic lipids capable of facilitating transport of biol. active agents or substances into cells, are disclosed. Thus, I (R1,R4 = oleyl; R2,R5 = Me2N(CH2)3; R3,R6 = Me; A1,A2 = CH2; L = (CH2)4; X = I) (II) is prepared by reaction of bis-1,4-oleyl-1,4-butanediamine with acrylonitrile followed by reduction of nitrile to amine and quaternization of amine with Me iodide.

II shows an activity of 37.8 ng/Pgal/cm2 in DNA delivery. Formulations containing I are given.
ACCESSION NUMBER: 2000:335366 CAPLUS
DOCUMENT NUMBER: 132:334312
TITLE: synthesis and activity of transfection reagents for transport of biol. active agents or substances into cells
INVENTOR(S): Chu, Yongliang; Masoud, Malek; Gebeyehu, Gulilat
PATENT ASSIGNEE(S): Life Technologies, Inc., USA
SOURCE: PCT Int. Appl., 130 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2000027795	A1	20000518	WO 1999-US26825	19991112
W: AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, DE, DK, DM, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
RW: GH, GM, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GN, GW, ML, MR, NE, SN, TD, TG				
CA 2350882	AA	20000518	CA 1999-2350882	19991112

L37 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2005 ACS ON STN (Continued)
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(synthesis and activity of transfection reagents for transport of biol. active agents or substances into cells)
RN 268539-49-7 CAPLUS
CN 2,6,11,15-Tetraazahexadecanediamide, 4,13-dihydroxy-6,11-di-(9Z)-9-octadecenyl- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

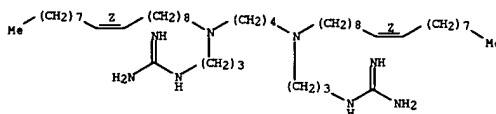


REFERENCE COUNT: 13 THERE ARE 13 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L37 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2005 ACS ON STN (Continued)
EP 1129064 A1 20010905 EP 1999-971794 19991112
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO
JP 2002529439 T2 20020910 JP 2000-580975 19991112
NZ 512244 A 20031219 NZ 1999-512244 19991112
AU 772847 B2 20040506 AU 2000-14776 19991112
PRIORITY APPLN. INFO.: US 1998-108117P P 19981112
WO 1999-US26825 W 19991112

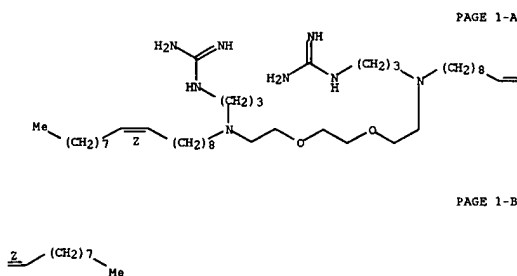
OTHER SOURCE(S): MARPAT 132:334312
IT 268539-59-9P 268539-60-2P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(synthesis and activity of transfection reagents for transport of biol. active agents or substances into cells)
RN 268539-59-9 CAPLUS
CN 2,6,11,15-Tetraazahexadecanediamide, 6,11-di-(9Z)-9-octadecenyl- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 268539-60-2 CAPLUS
CN 9,12-Dioxo-2,6,15,19-tetraazaeicosanediamide, 6,15-di-(9Z)-9-octadecenyl- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



IT 268539-49-7P

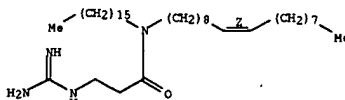
L37 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2005 ACS ON STN
AB Cationic lipids, RINCO(CH2)mNHCOCH(NHR2)(CH2)nNHR3 [R, R1 = alkyl; R2, R3 = H, acyl, alkyl, carboxamide, aryl; m = 1-3; n = 2-5], were prepared as permeability enhancer mols. to increase membrane permeability and facilitate cellular uptake of neg. charged comds. Thus, JA 59312, i.e. (Z)-H2NC(=NH)NH(CH2)2CH(NH2)CONHCH2CON(CH2)8CH(CH2)7Me (CH2)15Me, was prepared in a multistep synthetic scheme starting from 1-bromohexadecane, oleylamine, N-(benzyloxycarbonyl)-glycine N-hydroxysuccinimide ester, 2,4-bis[(tert-butoxycarbonyl)amino]butyric acid, and 1H-pyrazole-1-carboxamide hydrochloride. The prepared comds. were tested for cellular transport and cytotoxicity.

ACCESSION NUMBER: 1999:96203 CAPLUS
DOCUMENT NUMBER: 130:168544
TITLE: Preparation of novel compositions for the delivery of negatively charged molecules
INVENTOR(S): Beigelman, Leonid; Matulic-Adamic, Jasenka; Karpeisky, Alex; Haeblerli, Peter; Sweedler, David; Reynolds, Mark; Chaudhary, Wilab; Min, John
PATENT ASSIGNEE(S): Ribosome Pharmaceuticals, Incorporated, USA
SOURCE: PCT Int. Appl., 90 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: English
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9905094 A1		19990204	WO 1998-US15129	19980722
W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GE, GH, HU, IL, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
RW: AT, BE, BF, BJ, CF, CG, CH, CI, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG				
PRIORITY APPLN. INFO.: US 1997-53517 19970723 US 1998-72967 19980129				

OTHER SOURCE(S): MARPAT 130:168544
IT 220458-60-6P AK 524-73
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
(preparation of novel comds. for the delivery of neg. charged mols.)
RN 220458-60-6 CAPLUS
CN Propanamide, 3-[(aminomimomethyl)amino]-N-hexadecyl-N-(9Z)-9-octadecenyl- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



REFERENCE COUNT: 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS

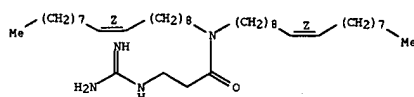
L37 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2005 ACS on STN
AB The title compds. R1R2NC(O)AX [R1, R2 = C10 - C26 hydrocarbyl; A = hydrocarbylene (further details on said hydrocarbylene are given); X = NHC((NR3)NR4, etc.); R3, R4 = hydrocarbyl, etc.; a proviso is given] are prepared in an in vivo gene transfer test, the transfection efficiency obtained with 2-quinidino-N,N-di-octadeca-9-enylpropionamide was greater than that achieved with Dotma.

ACCESSION NUMBER: 1998:379115 CAPLUS
DOCUMENT NUMBER: 129:81526
TITLE: Preparation of cationic lipids as materials for liposomes for gene transfer
INVENTOR(S): Belloni, Paula Nanette; Hirschfeld, Donald Roy; Rink, John Otto; Nester, John Joseph; Peltz, Gary Allen
PATENT ASSIGNEE(S): F. Hoffmann-La Roche A.-G., Switz.
SOURCE: Jpn. Kokai Tokkyo Koho, 29 pp.
CODEN: JXOXAF
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
JP 10152461	A2	19980609	JP 1997-285925	19971020
CA 2217550	AA	19980422	CA 1997-2217550	19971007
EP 846680	A1	19980610	EP 1997-117934	19971016
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO				
US 6034137	A	20000307	US 1997-954428	19971020
CN 1180697	A	19980506	CN 1997-121514	19971021
CN 1068585	B	20010718		
BR 9705117	A	19980915	BR 1997-5117	19971022
PRIORITY APPLN. INFO.: US 1996-29581P P 19961022 US 1997-49922P P 19970618				

OTHER SOURCE(S): MARPAT 129:81526
IT 209396-70-3P 209396-72-5P 209396-85-0P
209396-86-1P 209396-88-3P 209396-92-9P
209396-94-1P 209396-98-5P 209397-00-2P
209397-06-8P 209397-07-9P 209397-09-1P
209397-10-4P 209397-11-5P 209397-12-6P
RL: SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of cationic lipids as materials for liposomes)
RN 209396-70-3 CAPLUS
CN Propanamide, 3-[(aminoiminomethyl)amino]-N,N-di-(9Z)-9-octadecenyl-, monohydrochloride (9CI) (CA INDEX NAME)

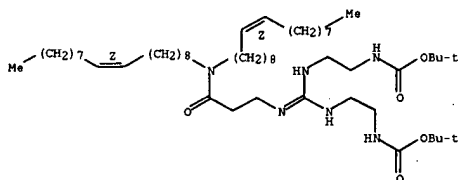
Double bond geometry as shown.



• HCl

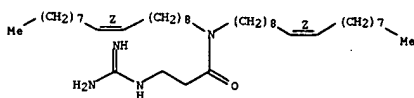
RN 209396-72-5 CAPLUS
CN 2,5,7,10-Tetraazadec-5-enedioic acid, 6-[[3-[di-(9Z)-9-octadecenylamino]-3-oxopropyl]amino]-, bis(1,1-dimethylethyl) ester (9CI) (CA INDEX NAME)

Double bond geometry as shown.



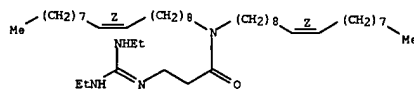
RN 209396-85-0 CAPLUS
CN Propanamide, 3-[(aminoiminomethyl)amino]-N,N-di-(9Z)-9-octadecenyl- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



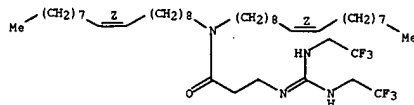
RN 209396-86-1 CAPLUS
CN Propanamide, 3-[[bis(ethylamino)methylene]amino]-N,N-di-(9Z)-9-octadecenyl- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



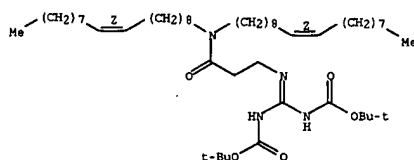
RN 209396-88-3 CAPLUS
CN Propanamide, N,N-di-(9Z)-9-octadecenyl-3-[[[(2,2,2-trifluoroethyl)amino]((2,2,2-trifluoroethyl)imino)methyl]amino]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



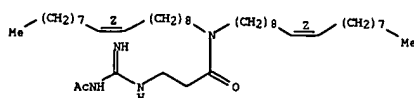
RN 209396-92-9 CAPLUS
CN Carbamic acid, [[3-[di-(9Z)-9-octadecenylamino]-3-oxopropyl]carbonimidoyl]bis-, bis(1,1-dimethylethyl) ester (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 209396-94-1 CAPLUS
CN Propanamide, 3-[[[(acetylamino)iminomethyl]amino]-N,N-di-(9Z)-9-octadecenyl- (9CI) (CA INDEX NAME)

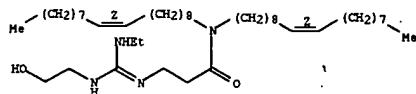
Double bond geometry as shown.



L37 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2005 ACS on STN (Continued)

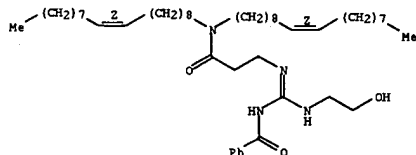
RN 209396-98-5 CAPLUS
CN Propanamide, 3-[[[ethylamino] (2-hydroxyethyl)amino]methylene]amino]-N,N-di-(9Z)-9-octadecenyl- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



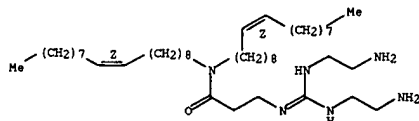
RN 209397-00-2 CAPLUS
CN Benzamide, N-[[[3-[di-(9Z)-9-octadecenylamino]-3-oxopropyl]amino] (2-hydroxyethyl)amino]methylene]- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 209397-06-8 CAPLUS
CN Propanamide, 3-[[[(2-aminoethyl)amino] (2-aminoethyl)imino]methyl]amino]-N,N-di-(9Z)-9-octadecenyl- (9CI) (CA INDEX NAME)

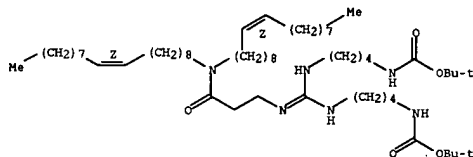
Double bond geometry as shown.



RN 209397-07-9 CAPLUS
CN Propanamide, 3-[[[bis(butylamino)methylene]amino]-N,N-di-(9Z)-9-octadecenyl- (9CI) (CA INDEX NAME)

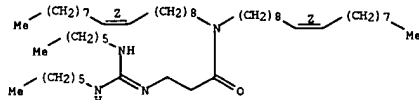
Double bond geometry as shown.

L37 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2005 ACS on STN (Continued)

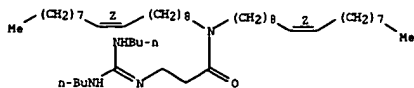


RN 209397-12-6 CAPLUS
CN Propanamide, 3-[[[bis(hexylamino)methylene]amino]-N,N-di-(9Z)-9-octadecenyl- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

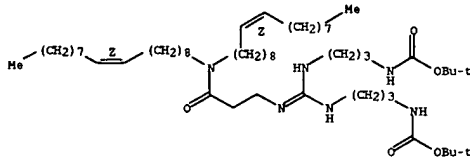


L37 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2005 ACS on STN (Continued)



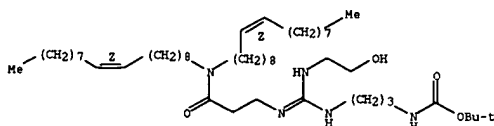
RN 209397-09-1 CAPLUS
CN 2,6,8,12-Tetraazatridec-6-enedioic acid, 7-[[[3-[di-(9Z)-9-octadecenylamino]-3-oxopropyl]amino]-, bis(1,1-dimethylethyl) ester (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 209397-10-4 CAPLUS
CN 2,6,8,12-Tetraazatridec-6,21-dienoic acid, 7-[[[3-[di-(9Z)-9-octadecenyl-11-oxo-, 1,1-dimethylethyl ester, (21E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.



RN 209397-11-5 CAPLUS
CN 2,7,9,14-Tetraazapentadec-7-enedioic acid, 8-[[[3-[di-(9Z)-9-octadecenylamino]-3-oxopropyl]amino]-, bis(1,1-dimethylethyl) ester (9CI) (CA INDEX NAME)

Double bond geometry as shown.

=> logoff y

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

18.42

1411.94

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL

ENTRY

SESSION

CA SUBSCRIBER PRICE

-2.19

-13.14

STN INTERNATIONAL LOGOFF AT 15:30:41 ON 25 APR 2005